# PRODUCT INFORMATION



## (R)-(+)-Linoleyl-1'-Hydroxy-2'-Propylamide

Item No. 9001233

CAS Registry No.: 220556-74-1

N-[(1R)-2-hydroxy-1-methylethyl]-Formal Name:

9Z,12Z-octadecadienamide

MF:  $C_{21}H_{39}NO_{2}$ 337.5 FW: **Purity:** ≥98%

Supplied as: A solution in ethanol

Storage: -20°C Stability: ≥2 years

Information represents the product specifications. Batch specific analytical results are provided on each certificate of analysis.

### **Laboratory Procedures**

(R)-(+)-Linoleyl-1'-hydroxy-2'-propylamide is supplied as a solution in ethanol. To change the solvent, simply evaporate the ethanol under a gentle stream of nitrogen and immediately add the solvent of choice. Solvents such as ethanol, DMSO, and dimethyl formamide purged with an inert gas can be used. The solubility of (R)-(+)-linoleyl-1'-hydroxy-2'-propylamide in these solvents is approximately 20, 5, and 11 mg/ml, respectively.

(R)-(+)-Linoleyl-1'-hydroxy-2'-propylamide is sparingly soluble in aqueous buffers. For maximum solubility in aqueous buffers, the ethanolic solution of (R)-(+)-linoleyl-1'-hydroxy-2'-propylamide should be diluted with the aqueous buffer of choice. (R)-(+)-Linoleyl-1'-hydroxy-2'-propylamide has a solubility of approximately 11 mg/ml in a 1:2 solution of ethanol:PBS (pH 7.2) using this method. We do not recommend storing the aqueous solution for more than one day.

#### Description

N-Acyl ethanolamines (NAEs) have diverse biological actions that are strongly affected by the associated acyl group. Linoleoyl ethanolamide (LOEA) has potential signaling roles in aging and neurological functioning. 1,2 LOEA has a weak affinity for cannabinoid (CB) receptors (K<sub>i</sub> = 10, 25 μM for CB<sub>1</sub>, CB<sub>2</sub>, respectively).<sup>3</sup> Although hydrolized by fatty acid amide hydrolase (FAAH; K<sub>i</sub> = 9 μM) it also inhibits FAAH and inhibits voltage-gated K<sup>+</sup> channels.<sup>3-6</sup> (R)-(+)-Linoleyl-1'-hydroxy-2'-propylamide is a homolog of LOEA, characterized by the addition of an (R)-α-methyl group at the methylene carbon adjacent to the amide nitrogen. A similar modification of arachidonoyl ethanolamide (Item No. 90050) to produce R-1 methanandamide (Item No. 90070) imparts higher affinity for the CB receptor as well as improved metabolic stability. The physiological actions of this compound have not been evaluated.

#### References

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- Lin, S., Khanolkar, A.D., Fan, P., et al. J. Med. Chem. 41, 5353-5361 (1998).
- 4. Maccarrone, M., van der Stelt, M., Rossi, A., et al. J. Biol. Chem. 273, 32332-32339 (1998).
- 5. Bisogno, T., Maurelli, S., Melck, D., et al. J. Biol. Chem. 272, 3315-3323 (1997).
- 6. Poling, J.S., Rogawski, M.A., Salem, N., Jr., et al. Neuropharmacology 35(7), 983-991 (1996).
- 7. Abadji, V., Lin, S., Taha, G., et al. J. Med. Chem. 37, 1889-1893 (1994).

WARNING
THIS PRODUCT IS FOR RESEARCH ONLY - NOT FOR HUMAN OR VETERINARY DIAGNOSTIC OR THERAPEUTIC USE.

This material should be considered hazardous until further information becomes available. Do not ingest, inhale, get in eyes, on skin, or on clothing. Wash thoroughly after handling. Before use, the user must review the complete Safety Data Sheet, which has been sent via email to your institution.

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